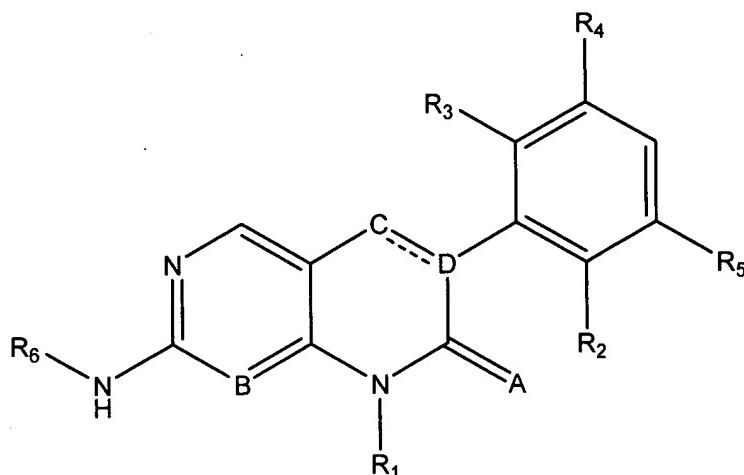


Amendments to the Claims:

This listing of claims will replace all prior versions and listings of claims in the application:

Listing of Claims:

1. (Currently Amended) A compound of the formula:



wherein:

A is O, NH₂, NH(C₁-C₆ alkyl), N(C₁-C₆ alkyl)₂, or -NHC(O)NHR₁₂;

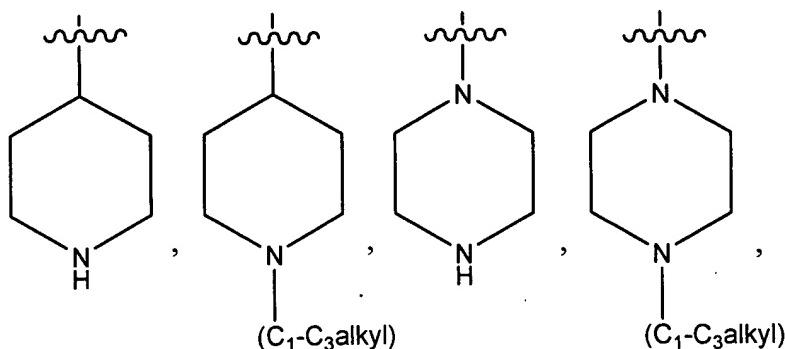
R₁₂ is C₁-C₆ straight or branched chain alkyl, or -(CH₂)_n-C₃-C₈ cycloalkyl ring; n is an integer of from 1 to 3;

~~B, C and D are independently selected from CH or N, with the proviso that C and D are not both N; B is N, C is CH, D is N.~~

R₁ is selected from the group of C₁-C₆ straight or branched chain alkyl, optionally substituted by -COOH, or;

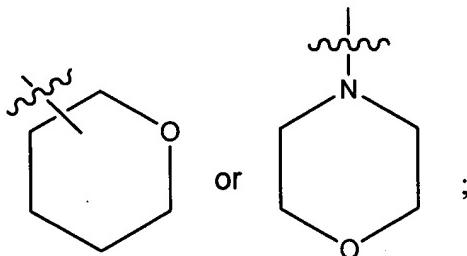
a) a phenyl, benzyl or C₃-C₈ cycloalkyl ring, or -CH₂-C₃-C₈ cycloalkyl ring, with the phenyl, benzyl or cycloalkyl rings being optionally substituted by 1 or 2 COOH or -CH₂-COOH groups; or

b) a piperidine or piperazine moiety selected from group of:



the rings of the piperidine or piperazine moieties being optionally substituted by 1 or 2 COOH or -CH₂-COOH groups; or

c) a tetrahydropyran or morpholine moiety of the formulae:



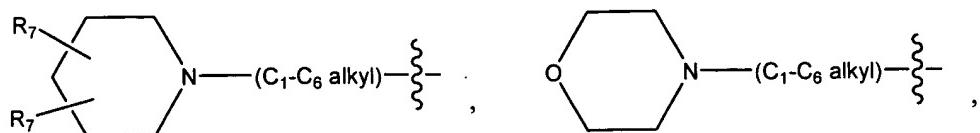
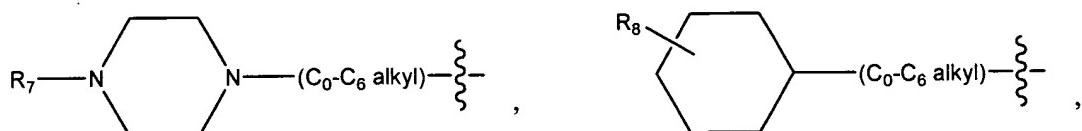
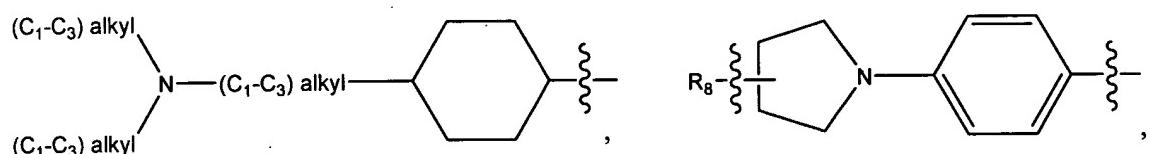
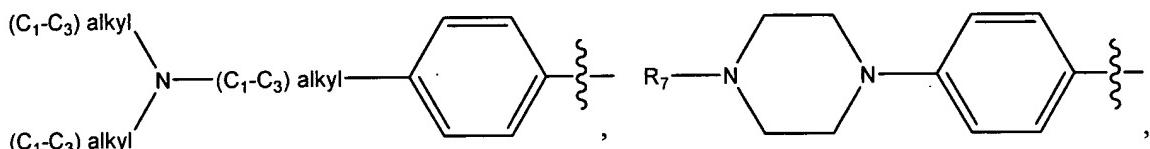
R₂ is H, Cl or F;

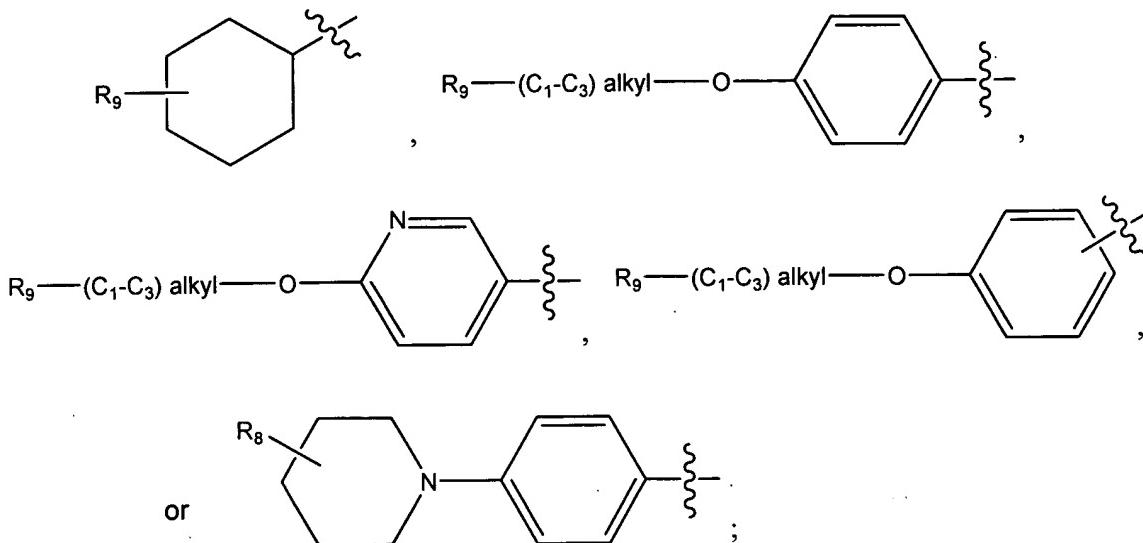
R₃ is H, Cl or F, with the proviso that at least one of R₂ or R₃ is F;

R₄ is H, OH, -OCH₃, or -OCH₂CH₃, with the proviso that, if R₄ is H, R₂ and R₃ are not H;

R₅ is -OCH₃, or -OCH₂CH₃;

R₆ is selected from the group of H, -(C₁-C₅ alkyl)-NH₂, -(C₁-C₅ alkyl)-NH-(C₁-C₃ alkyl)-R₁₁, -(C₁-C₅ alkyl)-N-(C₁-C₃ alkyl-R₁₁)₂, -O-(C₁-C₅ alkyl)-NH₂, -O-(C₁-C₅ alkyl)-NH-(C₁-C₃ alkyl)-R₁₁, -O-(C₁-C₅ alkyl)-N-(C₁-C₃ alkyl-R₁₁)₂, -CH(CH₂OH)₂, -(C₁-C₃ alkyl)(CH₂OH)₂, -(C₁-C₃ alkyl)-O-(C₁-C₃ alkyl)-R₁₁, -(C₁-C₃ alkyl)-O-(C₁-C₃ alkyl)-NH₂, -(C₁-C₃ alkyl)-O-(C₁-C₃ alkyl)-NH-(C₁-C₃ alkyl)-R₁₁, -(C₁-C₃ alkyl)-O-(C₁-C₃ alkyl)-N(C₁-C₃ alkyl-R₁₁)₂, phenyl substituted by one or two groups selected from NH₂, -N(C₁-C₃ alkyl), -N(C₁-C₃ alkyl)₂, CN or -(C₁-C₃ alkyl)-tetrazole, or C₁-C₆ alkyl,





with each of the alkyl chains of any group in this R_4 definition being optionally substituted by from 1 to 4 OH groups;

R_7 in each instance is independently selected from H, -NH₂, NH(C₁-C₃ alkyl), N(C₁-C₃ alkyl)₂, or C₁-C₃ alkyl;

R_8 is H, OH or C₁-C₃ alkyl;

~~9~~ R_9 is H, OH, -NH₂, NH(C₁-C₃ alkyl), or N(C₁-C₃ alkyl)₂;

R_{10} is H or C₁-C₃ alkyl;

R_{11} is H, CN, OH, NH₂, F, or CF₃,

or a pharmaceutically acceptable salt or ester form thereof.

2. (Currently amended) A compound of Claim 1 selected from the group of:

1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3-hydroxy-5-methoxy phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(S,S)-1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-[(5-hydroxymethyl-2-phenyl-[1,3]dioxolan-4-ylmethyl)-amino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

(S,S)-1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3,4-tr1Hydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(4hydroxy-cyclohexyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2-hydroxy-1-hydroxy-methyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

7-(3-Amino-2-hydroxy-propylamino)-1-cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;

1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-[2-(2-hydroxy-ethoxy)-ethylamino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Cyclopropyl-7-(4-diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or
7-(4-Diethylamino-butylamino)-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
or a pharmaceutically acceptable salt or ester form thereof.

3. (Currently amended) A compound of Claim 1 selected from the group of:

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-(2-hydroxy-1-hydroxymethyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-(4-hydroxy-cyclohexylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxyphenyl)-7-(2,3-dihydroxybutylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxypropyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or
7-(4-Amino-2,3-dihydroxy-butylamino)-1-cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-([S,S]-2,3,4-tr1Hydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-1-ethyl-7-[2-(2-hydroxyethoxy)-ethyl-amino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
7-(4-Amino-2,3-dihydroxy-butylamino)-3-(2,6-difluoro-3,5-dimethoxyphenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
7-{3-[Bis-(2-hydroxy-ethyl)-amino]-propylamino}-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or
1-Cyclopentyl-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxypropyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
or a pharmaceutically acceptable salt or ester form thereof.

4. (Currently amended) A compound of Claim 1 selected from the group of:

3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxy-butylamino)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
Ethyl-4-[3-(2,6-difluoro-3,5-dimethoxy-phenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-pyrimido[4,5-d]pyrimidin-1-yl]-cyclohexanecarboxylate;
4-[3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-

pyrimido[4,5-d]pyrimidin-1-yl]-cyclohexanecarboxylic acid;
7-Amino-3-(2,6-difluoro-3,5-dimethoxy-phenyl)-1-piperidin-4-yl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; compound with trifluoroacetic acid; or
1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(2-hydroxy-1-hydroxymethyl-ethylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Ethyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(4-hydroxy-butylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
(S)-1-Cyclopentyl-7-(2,3-dihydroxypropylamino)-3-(2-fluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Ethyl-3-(2-fluoro-3,5-dimethoxyphenyl)-7-(4-hydroxy-cyclohexylamino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
4-[3-(2,6-Difluoro-3,5-dimethoxyphenyl)-7-methylamino-2-oxo-3,4-dihydro-2H-pyrimido[4,5-d]pyrimidin-1-yl]-piperidine-1-carboxylic acid; or
3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-methylamino-1-piperidin-4-yl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
3-(2,6-Difluoro-3,5-dimethoxy-phenyl)-7-(2,3-dihydroxy-propylamino)-1-ethyl-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Cyclopentyl-7-(4-diethylamino-butylamino)-3-(3-ethoxy-2,6-difluoro-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
(R)-1-Cyclopentyl-7-(2,3-dihydroxy-propylamino)-3-(2-fluoro-3,5-dimethoxy-phenyl)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
1-Ethyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-[2-(2-hydroxy-ethoxy)-ethylamino]-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one; or
1-Cyclopentyl-3-(2-fluoro-3,5-dimethoxy-phenyl)-7-(*trans*-4-hydroxycyclohexyl-amino)-3,4-dihydro-1H-pyrimido[4,5-d]pyrimidin-2-one;
or a pharmaceutically acceptable salt or ester form thereof.

5. A pharmaceutical composition comprising a pharmaceutically effective amount of a compound of Claim 1 and a pharmaceutically acceptable carrier.

6-9. (Canceled)